

**Amendments to the Specification:**

On page 32, line 21, please replace Table IV with the following Table IV:

*TABLE IV. Coupling constants (*J*, Hz) of galactopyranose and glycerol residues of glycolipid derivatives and related reference compounds*

	BBGL-II	BBGL-II Ac <sub>3</sub>	BBGL-III	BBGL-III Ac <sub>4</sub>	MDGD	1,2-Dipamitin	Mc- $\alpha$ -D-galp	Mc- $\beta$ -D-galp
<b>Solvent</b>								
CDCl <sub>3</sub>	CDCl <sub>3</sub>	CDCl <sub>3</sub>	CDCl <sub>3</sub>	CDCl <sub>3</sub> : CD <sub>3</sub> OD (4:1 v/v)				
<i>J</i> <sub>1,2</sub> (Gal)	7.5	8.0	3.8	3.7	7.3	3.4	8.0	
<i>J</i> <sub>2,3</sub>	ND <sup>a</sup>	10.5	9.8	10.0	9.7	10.1	9.9	
<i>J</i> <sub>3,4</sub>	3.2 <sup>b</sup>	3.5	3.2	3.5	3.3	2.8	3.5	
<i>J</i> <sub>5,6</sub>	1.0 <sup>b</sup>	1.0	1.1	1.1	1.1	1.6	0.8	
<i>J</i> <sub>5,6a</sub>	6.3	6.7	5.1	6.4	6.5	6.8	7.9	
<i>J</i> <sub>5,6b</sub>	7.2	6.7	ND	7.0	5.4	5.5	4.4	
<i>J</i> <sub>6a,6b</sub>	11.1	11.2	11.5	11.2	11.6	11.7	11.7	
<i>J</i> <sub>C-I' H-I'</sub>	158.7	157.4	170.5	172.4	160.2	170.2	160.6	
<i>J</i> <sub>1,3a,b</sub> <sup>c</sup> (Gro)			11.9	11.8	12.1	11.9		
<i>J</i> <sub>1,3a,b,2"</sub>			4.1	4.1	3.2	4.5		
<i>J</i> <sub>1,3a,b,2"</sub>			5.9	6.1	6.7	5.7		
<i>J</i> <sub>2",3",3"</sub> <sub>a</sub>			4.8	4.3	5.4	4.8		
<i>J</i> <sub>2",3",3"</sub> <sub>b</sub>			6.2	5.2	6.0	5.2		
<i>J</i> <sub>3",a,3"</sub> <sub>b</sub>			10.9	11.2	10.9	12.2		

<sup>a</sup> Not determined.

<sup>b</sup> CD<sub>3</sub>)<sub>2</sub>CO:CDCl<sub>3</sub> (7:3 v/v) solution.